

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102418\
 Data File : VX005543.D
 Acq On : 24 Oct 2018 16:06
 Operator : JC/MD
 Sample : J5655-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S37-EB-2018-1

Quant Time: Oct 25 05:19:05 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102418W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 24 13:13:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	235724	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	328411	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	288711	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	134533	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	138247	48.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.08%	
35) Dibromofluoromethane	5.50	113	107467	48.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.66%	
50) Toluene-d8	8.71	98	385497	48.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.50%	
62) 4-Bromofluorobenzene	11.14	95	126414	41.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.78%	
Target Compounds						
16) Acetone	2.45	43	4805	3.160	ug/l	94
20) Methylene Chloride	2.86	84	46319	18.189	ug/l	97
25) 2-Butanone	4.70	43	6449	2.785	ug/l	99
29) Tetrahydrofuran	5.15	42	169482	112.477	ug/l	92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX102418\
Data File : VX005543.D
Acq On : 24 Oct 2018 16:06
Operator : JC/MD
Sample : J5655-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
S37-EB-2018-1

Quant Time: Oct 25 05:19:05 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X102418W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 24 13:13:34 2018
Response via : Initial Calibration

